

Reduction of local velocity spreads by linear potentials

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We study the spreading of the wave function of a Bose-Einstein condensate accelerated by a constant force both in the absence and in the presence of atom-atom interactions. We show that, despite the initial velocity dispersion, the local velocity dispersion defined at a given position downward can reach ultralow values and be used to probe very narrow energetic structures. We explain how one can define quantum mechanically and without ambiguities the different velocity moments at a given position by extension of their classical counterparts. We provide a common theoretical framework for interacting and non-interacting regimes based on the Wigner transform of the initial wave function that encapsulates the dynamics in a scaling parameter. In the absence of interaction, our approach is exact. Using a numerical simulation of the 1D Gross-Pitaevskii equation, we provide the range of validity of our scaling approach and find a very good agreement in the Thomas-Fermi regime. We apply this approach to the study of the scattering of a matter wave packet on a double barrier potential. We show that a Fabry-Perot resonance in such a cavity with an energy width below the pK range can be probed in this manner. We show that our approach can be readily transposed to a large class of many-body quantum systems that exhibit self-similar dynamics.

I. INTRODUCTION

Narrowing down the velocity spread of a packet of particles or a beam is important in many techniques and experiments: to induce, control or measure energy-dependent phenomena, such as sharp resonances or energy thresholds; to probe, by scattering, surfaces or interaction potentials with high energy resolution; or to improve interferometers, spectroscopic studies, and metrological devices such as atomic clocks. A recent application example is the study of very low-energy molecular collisions thanks to an accurate velocity control of merged supersonic molecular beams [1].

Frequently the particles are produced at a preparation chamber and interact with a material target or field elsewhere after a free or guided flight. This means that the velocity spread of interest is the one of the particles as they arrive at the interaction region. The standard (global) velocity spread of the ensemble of particles, irrespective of their location, is only relevant inasmuch as it affects the local velocity spread, which, as we shall see, may differ substantially from the global one.

In this paper we propose and characterize a method to achieve a narrow local velocity distribution for one dimensional motion. Our main motivation is to set high resolution collision experiments between ultracold atoms and a well localized obstacle, but the principles involved are applicable beyond that goal. Local velocities and their statistics may be defined in classical mechanics since trajectories carry simultaneous position and velocity information, but in quantum mechanics position and momentum do not commute. This does not mean though that a quantum local velocity is a meaningless concept

[2]. A detailed analysis of specific detectors to perform a local velocity measurement would be necessary in order to determine the exact operators (i.e. the specific quantization among the many possible) and/or Positive Operator Valued Measures (POVMs) involved. This operational approach is out of the scope of this paper. Instead we shall adopt a simple phase-space Wigner-function description of the dynamics and the corresponding Weyl rule of quantization to specify the local quantum observables starting from classical expressions, so that we shall treat formally quantum and classical systems alike in phase-space. In any case we shall assume conditions in which the differences among different quantization rules are negligible.

A well-known narrowing of the local velocity distribution occurs for free motion. If a cloud of classical (non-interacting) particles with an initial phase-space density without position-momentum correlations and zero average position and velocity is let to expand freely, faster particles will advance beyond the slow ones, so that at a distant observation point the particles that arrive *at a specific instant of time* have well defined velocities. Thus this type of free-motion induced velocity narrowing only applies to the instantaneous local velocity distribution. The initial spread for all particles irrespective of their location is actually constant in time, and implies that the dominant, instantaneous (mean) velocity at the observation point changes with time. In other words, the instantaneous local velocity spread is small, but the spread in time of the instantaneous local mean velocities may be large. This limits severely the applicability and usefulness of free-motion narrowing. If we are interested in a specific velocity range, many, even most of the particles, i.e., those that do not arrive at the “right” time interval

to match our requirements, should be discarded. Instead, we propose a mechanism so that the velocities of all particles irrespective of their arrival time is sharply defined. It only requires a constant force, which for neutral cold atoms may be implemented by using gravity or a magnetic field gradient on magnetically polarized atoms.

We shall study the experimental situation depicted in Fig. 1. In Secs. II and III, the general formalism to characterize the velocity distribution at the obstacle location x is worked out. In Sec. IV, we will apply it to noninteracting particles, and in Sec. V to interacting particles in the Thomas-Fermi regime. In Sec. VI, a full numerical study of the scattering on a double well potential from a carefully outcoupled matterwave is carried out. It shows the possibility to investigate matter wave Fabry-Perot cavities.

II. GENERAL FORMALISM

The statistical analysis we perform is based on a double average, over velocity and time, at the observation location. To that end we follow, interchanging the roles of time and position, the double averaging performed in [2] over position and momentum at fixed time. The definition of the local instantaneous averages over the arriving velocities could in principle be carried out in several ways with different physical implications. For a phase-space distribution $W(q, p; t)$ normalized to one at any time t when integrating over positions q and momenta p , a local, instantaneous average at $q = x$ and time t may be defined as

$$v_x(t) = \frac{\int dp v W(x, p; t)}{\int dp W(x, p; t)}, \quad (1)$$

where $v = p/m$ and m is the mass of the particle. This average is known in hydrodynamical formulations as the “velocity field”, see e.g. [3, 4], and it becomes operationally meaningful if the velocities of the particles between x and $x + dx$ are probed instantaneously. The normalization factor in the denominator is just the particle density divided by the total number of atoms N . For many experiments and detectors the particles arrive according to a certain time-dependent flux and the relevant quantity is not the number of particles *present* in dx with a specific velocity, but rather the number of particles *arriving* (crossing x) with velocities between v and $v + dv$ in dt . This is given by $NvW(x, p, t)dvdt$ so that the local instantaneous mean velocity (and similarly for higher moments) is defined as

$$\bar{v}_x(t) = \frac{\int dp v^2 W(x, p; t)}{\int dp v W(x, p; t)}, \quad (2)$$

where the denominator is now the current density $J_x(t)$ or flux (per particle) instead of the density. This is in particular the average applicable for scattering experiments [5–8]. A basic assumption is that all particles arrive from

the left (all velocities are positive at x) so $J_x(t) \geq 0$ and $\int dt J_x(t) = 1$.

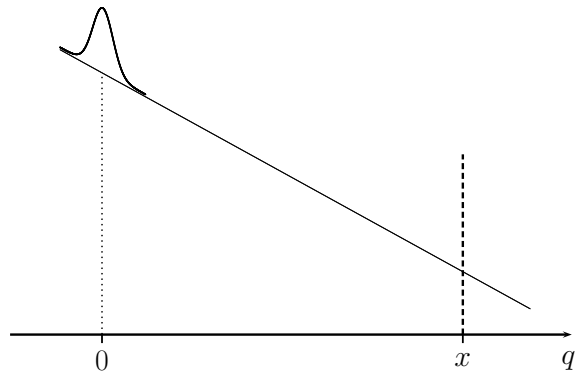


FIG. 1: A wave packet initially located at $q = 0$ and subjected to a constant force evolves towards an interacting region located at $q = x$.

Let us first assume a packet of classical non-interacting atoms described by a phase-space-density distribution $W(q, p; t)$ normalized to one. This can be translated into quantum mechanics by means of phase-space quasidistribution functions. For the problem at hand (a linear-in- q potential) the Wigner function is a natural choice, since its propagator for non-interacting particles is identical in classical and quantum mechanics [2].

The fraction of the number of atoms $d^2N_x^{(v)}(t)$ with a velocity between v and $v + dv$ that crosses the plane $q = x$ between t and $t + dt$ is

$$d^2N_x^{(v)}(t) = W(x, p; t) v dp dt \equiv P_x(v, t) dv dt. \quad (3)$$

The probability distribution $P_x(v, t)$ represents the probability density of particles in velocity space per unit time that crosses the plane x at t . We assume that all particles have a positive velocity when crossing x , so

$$\iint P_x(v, t) dv dt = \int J_x(t) dt = 1. \quad (4)$$

The quantity $J_x(t)$ corresponds to the local atomic flux divided by N ,

$$J_x(t) = \frac{dN_x(t)}{dt} = \int \frac{d^2N_x^{(v)}(t)}{dt} = \int W(x, p; t) v dp, \quad (5)$$

where N_x is the probability to find a particle at $q > x$. In order to define local instantaneous averages it is convenient to introduce the conditional probability density $P_x(v|t)$: $P_x(v|t)dv$ represents the probability to find the velocity between v and $v + dv$ for the particles that cross x between t and $t + dt$. It is normalized to one,

$$\int P_x(v|t) dv = 1. \quad (6)$$

According to Bayes law, the simple relation

$$P_x(v, t) = P_x(v|t) J_x(t) \quad (7)$$

holds. The instantaneous mean local velocity at position x and time t is given by

$$\bar{v}_x(t) = \int v P_x(v|t) dv. \quad (8)$$

Similarly, the instantaneous local mean square velocity reads

$$\overline{v_x^2}(t) = \int v^2 P_x(v|t) dv. \quad (9)$$

The mean local velocity is defined as an average over time of the instantaneous mean local velocity,

$$\begin{aligned} \langle v \rangle_x &= \iint v P_x(v, t) dv dt = \iint v^2 W(x, p; t) dp dt \\ &= \int \bar{v}_x(t) J_x(t) dt. \end{aligned} \quad (10)$$

Similarly, the local variance of the velocity reads

$$\begin{aligned} (\Delta v_x)^2 &\equiv \langle (v - \langle v \rangle_x)^2 \rangle_x = \iint (v - \langle v \rangle_x)^2 P_x(v, t) dv dt \\ &= \iint (v - \bar{v}_x + \bar{v}_x - \langle v \rangle_x)^2 P_x(v, t) dv dt \\ &= \iint [(v - \bar{v}_x)^2 + (\bar{v}_x - \langle v \rangle_x)^2] P_x(v, t) dv dt \\ &= \int \left[\int (v - \bar{v}_x)^2 P_x(v|t) dv + (\bar{v}_x - \langle v \rangle_x)^2 \right] J_x(t) dt \\ &= \int [\sigma_{v|x}^2(t) + D_{v|x}^2(t)] J_x(t) dt, \end{aligned} \quad (11)$$

where

$$\begin{aligned} \sigma_{v|x}^2(t) &\equiv \frac{1}{J_x(t)} \int (v - \bar{v}_x)^2 P_x(v, t) dv, \\ D_{v|x}^2(t) &\equiv (\bar{v}_x - \langle v \rangle_x)^2, \end{aligned} \quad (12)$$

are, respectively, the instantaneous local velocity variance and the squared deviation of the instantaneous local mean velocities with respect to the local average velocity. According to Eq. (11), the variance of the velocity at x is the sum of two terms: the time average of the instantaneous local variance and the variance of the instantaneous mean velocities.

Introducing the velocity moments

$$V_n = V_n(x, t) \equiv \int v^n W(x, p; t) dp, \quad (13)$$

we may rewrite the above quantities as

$$\begin{aligned} J_x(t) &= V_1(x, t), \\ \bar{v}_x(t) &= \frac{1}{J_x(t)} \int v^2 W(x, p, t) dp = \frac{V_2}{V_1}, \\ \overline{v_x^2}(t) &= \frac{1}{J_x(t)} \int v^3 W(x, p; t) dp = \frac{V_3}{V_1}, \\ \sigma_{v|x}^2(t) &= \frac{V_3}{V_1} - \left(\frac{V_2}{V_1} \right)^2, \\ \langle v \rangle_x &= \iint v^2 W(x, p, t) dp dt = \int dt V_2(x, t), \\ D_{v|x}^2(t) &= \left(\frac{V_2(x, t)}{V_1(x, t)} - \int V_2(x, t') dt' \right)^2, \end{aligned}$$

so computing the first three moments is enough to calculate the local variance of the velocity (11). The global velocity dispersion irrespective of the particle location is given by

$$\Delta v(t) = \sqrt{\int V_2(x, t) dx - \left(\int V_1(x, t) dx \right)^2}. \quad (14)$$

III. THE TIME-DEPENDENT WIGNER FUNCTION AND ITS VELOCITY MOMENTS

In the following, we study the dynamics of the ground state wave function of a 1D harmonic trap of angular frequency ω_0 after its sudden release on a linear potential. The scaling formalism developed hereafter is applied to a Bose-Einstein condensate (BEC) in the Thomas-Fermi regime [10–12] and to the Gaussian wave function for the non-interacting case. However, it is worth noticing that it can be applied to the class of many-body quantum systems that exhibit a self-similar dynamics [13, 38] as justified in Appendix A. In the presence of a constant acceleration field, γ , the time-dependent Wigner function, $W(x, p, t)$, can be obtained from the initial Wigner function, W_0 , through a proper scaling substitution (see Appendix A),

$$W(x, p, t) = W_0(X, P), \quad (15)$$

where we have used the following linear transformation

$$\begin{pmatrix} X \\ P \end{pmatrix} = \begin{pmatrix} 1/\alpha & 0 \\ -m\dot{\alpha} & \alpha \end{pmatrix} \begin{pmatrix} x - \eta \\ p - m\dot{\eta} \end{pmatrix}, \quad (16)$$

with α and η two time-dependent functions. The matrix of this transformation generates both a squeezing in momentum space correlated to a shearing in position space. η accounts for the center of the packet motion and obeys the Newton equation $\ddot{\eta} = \gamma$. The scaling factor α obeys a second order differential equation that depends on the system (see Appendix A):

$$\ddot{\alpha} = \frac{\omega_0^2}{\alpha^p}, \quad (17)$$

with $p = 3$ for a Gaussian wave function without interactions or a Tonks-Girardeau gas [14–16], and $p = 2$ for the wave function of a Bose-Einstein condensate in the Thomas-Fermi regime. The X and P variables are useful to derive the velocity moments in terms of their initial value $V_n^{(0)}(x)$:

$$V_n(x, t) = \int \left[\frac{P}{m\alpha} + u(X, t) \right]^n W_0(X, P) \frac{dP}{\alpha}, \quad (18)$$

with $u(X, t) = \dot{\eta} + \dot{\alpha}X$. We find

$$V_0(x, t) = V_0^{(0)}(X)/\alpha, \quad (19)$$

$$V_1(x, t) = uV_0^{(0)}(X)/\alpha, \quad (20)$$

$$V_2(x, t) = V_2^{(0)}(X)/\alpha^3 + u^2V_0^{(0)}(X)/\alpha, \quad (21)$$

$$V_3(x, t) = 3uV_2^{(0)}(X)/\alpha^3 + u^3V_0^{(0)}(X)/\alpha. \quad (22)$$

To obtain this result, we have explicitly used the symmetry property of the initial Wigner function: $W_0(x, p) = W_0(x, -p)$. The first three moments are therefore directly deduced from two moments of the initial distribution. Interestingly, the first two global velocity moments are readily obtained in a form that depends explicitly on the initial condition and the scaling parameters. We find $\langle v \rangle(t) = \dot{\eta}$ and the velocity dispersion

$$\Delta v^2(t) = \frac{\Delta v^2(0)}{\alpha^2} + \frac{2\omega_0^2}{p-1} (1 - \alpha^{1-p}) \Delta x^2(0), \quad (23)$$

using Eq. (17). With $p = 3$, we recover the fact that the velocity dispersion remains constant and equal to its initial value in the case of an interaction-free Gaussian wave packet. With $p = 2$ we obtain the explicit expression for the velocity dispersion at any time for a Bose-Einstein condensate in the Thomas-Fermi regime. In particular, its limit for long time is equal to $\Delta v(t \rightarrow \infty) = \sqrt{2}\omega_0\Delta x(0)$ in agreement with the Virial theorem [17, 18].

IV. APPLICATION TO A NONINTERACTING GAS

Let us set the initial phase-space distribution as an uncorrelated product of Gaussians,

$$W_0(x, p) = \frac{1}{2\pi} \frac{1}{\Delta x_0 \Delta p_0} e^{-x^2/2\Delta x_0^2} e^{-p^2/2\Delta p_0^2}, \quad (24)$$

obeying $\Delta x_0 \Delta p_0 = \hbar/2$. For a classical ensemble \hbar could be an arbitrary constant whereas in quantum mechanics this is Planck's constant/ 2π . Using $\sigma \equiv \Delta x_0$ we find

$$\begin{aligned} V_0^{(0)}(x) &= \frac{1}{\sqrt{2\pi}\sigma} e^{-x^2/2\sigma^2}, \\ V_2^{(0)}(x) &= \left(\frac{\hbar}{2m\sigma} \right)^2 V_0^{(0)}(x), \\ \alpha(t) &= \sqrt{1 + \omega_0^2 t^2}, \\ u(x, t) &= \gamma t + (x - \gamma t^2/2)\omega_0^2 t / (1 + \omega_0^2 t^2), \end{aligned} \quad (25)$$

with $\omega_0 = \hbar/(2m\sigma^2)$.

The quantities $\bar{v}_x(t)$, $\overline{v_x^2}(t)$ and $\sigma_{v|x}^2(t)$ involve ratios of the form $V_{n>1}(x, t)/V_1(x, t)$. Using Eqs. (20), (21) and (22), one realizes that these ratios depend only on $V_2^{(0)}(X)/V_0^{(0)}(X)$, which is a constant for a Gaussian Wigner function. In this way, we find

$$\bar{v}_x(t) = \frac{u^2 + (u^2 t^2 + \sigma^2)\omega_0^2}{u(1 + \omega_0^2 t^2)}, \quad (26)$$

$$\overline{v_x^2}(t) = u\bar{v}_x(t) + \frac{2\sigma^2\omega_0^2}{1 + \omega_0^2 t^2}. \quad (27)$$

The first two global velocity moments are

$$\langle v \rangle(t) = \int V_1(x, t) dx = \gamma t, \quad (28)$$

$$\langle v^2 \rangle(t) = \int V_2(x, t) dx = \frac{\hbar^2}{4m^2\sigma^2} + \gamma^2 t^2. \quad (29)$$

To have a negligible fraction of negative velocities we simply assume an observation point at $x \gg \sigma$ and $t > 0$ hereafter, in particular in all integrals. We now calculate the first velocity moments. They can be written in the form

$$V_j(x, t) := e^{-\Phi} f_j, \quad (30)$$

where the f_j are given in Appendix B and

$$\Phi = \frac{m^2\sigma^2(\gamma t^2 - 2x)^2}{8m^2\sigma^4 + 2\hbar^2 t^2} = -\frac{x^2(1 - \gamma t^2/2x)^2}{2\sigma^2(1 + \hbar^2 t^2/4m^2\sigma^4)} \quad (31)$$

is the phase that governs the behavior of the moments. In particular $\Phi = \Phi' = 0$ (the primes denote time derivatives) at $t_c = \sqrt{2x/\gamma}$, which is the time that a classical particle takes to reach x if it is initially at rest at the origin. The second derivative with respect to time at that instant is

$$\Phi''(t_c) = \frac{16gm^2\sigma^2 x}{8m^2\sigma^4 + 4\hbar^2 x/\gamma} \quad (32)$$

which, keeping x/γ constant, i.e., for a fixed t_c , grows with x . In other words, the moments become narrower functions of t .

We may now calculate the local variance as

$$(\Delta v_x)^2 = \int_0^\infty dt V_3 - \left[\int_0^\infty dt V_2 \right]^2. \quad (33)$$

Laplace's (or steepest descent) approximation gives for the integrals $\int e^{-\Phi} f dt \sim e^{-\Phi(t_c)} f(t_c) \sqrt{2\pi/\Phi''(t_c)}$ and $(\Delta v_x)^2 \sim 0$ in sharp contrast with the fixed-time (and time independent) global velocity variance $\langle v^2 \rangle(t) - (\langle v \rangle(t))^2 = \hbar^2/(4m^2\sigma^2)$ calculated with the Wigner function, see Eqs. (28) and (29). In Fig. 2 we show the two components that, when integrated, make $(\Delta v_x)^2$, see Eq. (11): In our configuration the local velocity variance

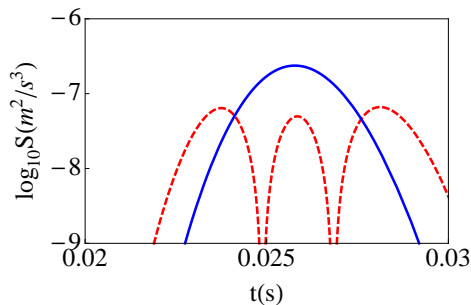


FIG. 2: (Color online) Components of the integrand for the local variance Δv_x^2 : $S = \sigma_{v|x}^2(t) J_x(t)$ (dashed red line), and $S = D_{v|x}^2(t) J_x(t)$ (solid blue line). Parameters: $\gamma = 0.6 \text{ m/s}^2$, $x = 201 \text{ } \mu\text{m}$, mass of rubidium-87. The initial state is the ground state of a harmonic trap with angular frequency $\omega_0 = 2\pi \times 150 \text{ Hz}$. The passage of a classical particle released from the origin at $t = 0$ is at $t_c = 0.026 \text{ s}$.

is dominated by the squared deviation of the local instantaneous mean. The parameters are chosen as specified in the caption for realistic conditions on initial velocity width, acceleration, and observation distance. For these parameters the local variance $(\Delta v_x)^2(0.5\text{mm}) = 8.22 \times 10^{-10} \text{ (m/s)}^2$, is already more than two orders of magnitude smaller than the global variance $(\Delta v)^2 = 3.4 \times 10^{-7} \text{ (m/s)}^2$.

The two variances (global and local) are quite different because of their distinct physical content. $(\Delta v)^2$ is independent of x , γ or t and corresponds to measuring the velocities of all the particles of the statistical ensemble, wherever they are, at a given instant. Since all particles are equally accelerated their initial velocity differences remain the same at any time. $(\Delta v_x)^2$ results instead from local velocity measurements weighted by $P_x(v, t)$. This distribution can be narrowly peaked when the velocities at x are predominantly due to the common effect of gx rather than to the effect of the initial spread of the initial conditions. Specifically we have that, at position x , the velocity of a particle that at time 0 was at x_0 with velocity v_0 , is

$$v_x = \sqrt{\gamma(x - x_0) + v_0^2}, \quad (34)$$

i.e., any difference due to the the initial conditions is washed out for a large $x\gamma$, which leads to velocity uniformity. The conditions $x \gg \sigma$ and $x\gamma \gg \sigma_v^2$, where σ_v^2 is the initial variance of the velocity, guarantee the dominance of the first term and lead to a local velocity spread suppression. Note that, when σ and σ_v are related by the minimum uncertainty product relation, i.e., $\sigma_v = \hbar/(2m\sigma)$, very small or very big σ values require a distant x to wash out the differences in the arriving velocities. Indeed, all previous results can be translated formally into quantum mechanics by interpreting W as the Wigner function and the dynamical variables according to Weyl's quantization rule. In the Weyl-Wigner framework the dynamical variable $\delta(x - q)v$ corresponds to the

flux operator $\hat{J}_x = (1/2)(\hat{v}\delta(x - \hat{q}) + \delta(x - \hat{q})\hat{v})$ and, more generally, $v^n \delta(x - q)$ to hermitian operators with matrix elements

$$\langle p | O^{(n)} | p' \rangle = \left(\frac{p + p'}{2m} \right)^n \langle p | x \rangle \langle x | p' \rangle \quad (35)$$

in momentum space. As other quantization rules are possible, the above results, which are based on one of them, may be questioned because of their non-uniqueness. However, actual measurements typically imply an averaging that tends to smooth the differences among the different rules. Moreover the conditions of the experiment discussed are such that at each instant the local quantum field is essentially monochromatic. This again washes out differences among quantization rules: as interferences do not play any major role for our setting and the effects of the non-commutativity of position and momentum are negligible, a classical-like interpretation of the results is justified. This is in fact the usual approach for most time-of-flight experiments.

V. APPLICATION TO A BOSE-EINSTEIN CONDENSATE IN THE THOMAS-FERMI REGIME

Unlike the Gaussian case, the Wigner function of the wave function associated with a Bose-Einstein condensate in the Thomas-Fermi regime cannot be worked out analytically. In this section, we compare numerics with the scaling approach to obtain the different velocity moments accurately. The parabolic profile of the condensate in the Thomas-Fermi regime has been used to deduce Eqs. (15) and (17) (see Appendix A) that gives the time evolution of the scaling parameter. However, a numerical study is necessary to validate the determination of the velocity moments with the scaling approach since the parabolic profile does not encapsulate the smooth edge of the wavefunction [19]. We show in the following that the knowledge of the initial Wigner function computed numerically is sufficient to obtain all required information on the dynamics and in particular on the different velocity moments with a high accuracy.

The initial wave function is found by evolving the 1D time-dependent Gross-Pitaevskii (GP) equation in imaginary time using a split-step Fourier method [20]. In practice, we use an initial Gaussian trial wavefunction with a mean quadratic size evaluated from the expected parabola shape and we iterate until the final wavefunction is stationary with respect to real-time evolution. The mean-field wave function corresponds to that of a Bose-Einstein condensate with N rubidium-87 atoms (scattering length $a_{sc} = 5 \text{ nm}$) held in a harmonic trap of angular frequency $\omega_0 = 2\pi \times 150 \text{ Hz}$. The Thomas-Fermi regime is reached when $\chi = Na_{sc}/a_0 \gg 1$ where $a_0 = (\hbar/m\omega_0)^{1/2}$ is the length scale associated with the harmonic confinement ($\chi = 1$ corresponds to $N \approx 176$ atoms).

In the following, we consider the time evolution of the wave function resulting from an abrupt release of the condensate on the slope ($V(x) = m\gamma x$, $\gamma = 0.6$ m/s²), the case of a progressive outcoupling is addressed in Sec. VI. We compare the results obtained by a full integration of the Gross-Pitaevskii equation with those resulting from the scaling approach (Eqs. (23) and (17)). In the course of the propagation, the interaction energy is converted into kinetic energy as illustrated in Fig. 3, and more precisely it is responsible for the increase of velocity dispersion. After 5 ms of propagation for a BEC initially in the Thomas-Fermi regime, already 80 % of the interaction energy has been transferred. Figure 4 shows the evolution of the velocity dispersion depending on the number of atoms. The impressive agreement demonstrates the efficiency of the scaling approach.

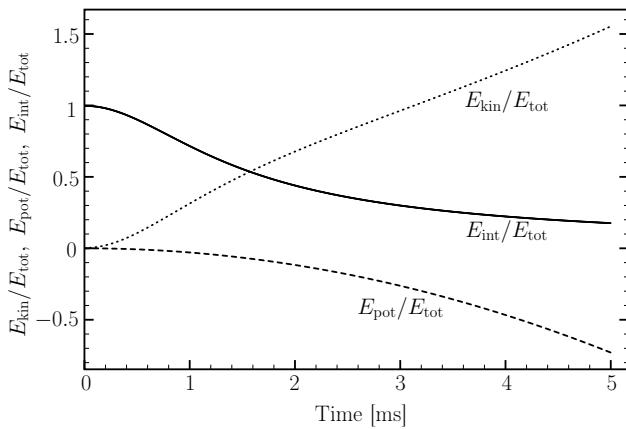


FIG. 3: Evolution of the kinetic (E_{kin} , dotted line), potential (E_{pot} , dashed line) and interaction energy (E_{int} , solid line) of a Bose-Einstein condensate as a function of time during its propagation over the slope (energies are normalized to the total energy $E_{\text{tot}} = E_{\text{kin}} + E_{\text{pot}} + E_{\text{int}}$). Parameters: the initial wave function corresponds to that of a Bose-Einstein condensate of rubidium-87 in the Thomas-Fermi regime with $N = 17600$ atoms initially at equilibrium in an harmonic potential of angular frequency $\omega_0 = 2\pi \times 150$ Hz. At time $t = 0$, the trapping potential is suddenly removed and the packet experiences the acceleration ($\gamma = 0.6$ m/s²) due to the linear potential.

The numerical integration of the GP equation with split-Fourier method is quite efficient even though the time step should be smaller for large-size wavefunction (large N). However, the computation of the Wigner function turns out to be rapidly cumbersome since the required grid in phase-space has a size that should increase with time to guarantee a good accuracy. We now compare the first two non-zero velocity moments $V_0(x, t)$ and $V_2(x, t)$ obtained after 5 ms of time evolution on the slope from the full numerical procedure of the dynamics (integration of the Gross-Pitaevskii equation + calculation of the Wigner function at time $t = 5$ ms) with the result of the scaling approach for which we use only the computing of the initial Wigner function combined with the

time evolution of the scaling parameter (Eqs. (16) and (17)). Results are summarized in Fig. 5 for $N = 35200$ atoms.

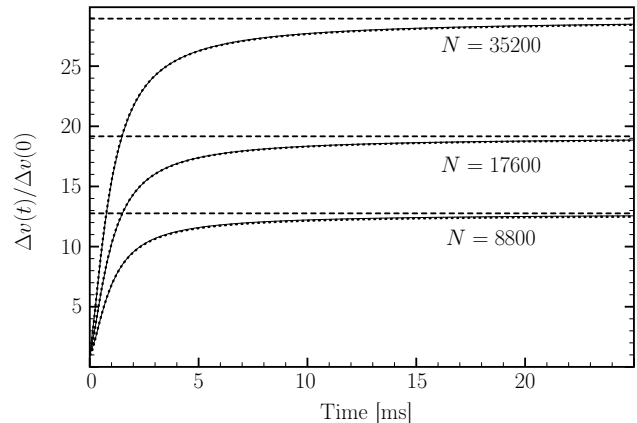


FIG. 4: Time evolution of the global velocity dispersion $\Delta v(t)/\Delta v(0)$ of a Bose-Einstein condensate for three different values of the atom number $N = 8800, 17600, 35200$. Solid lines: theoretical prediction using the scaling approach (see Eq. 23), dotted lines numerical integration of the Gross-Pitaevskii equation, dashed line: theoretical asymptote at long time.

To compare quantitatively the velocity moments, we define the relative error after 5 ms of propagation:

$$\varepsilon(N) = \left| 1 - \frac{\Delta_2 v(t = 5 \text{ ms})}{\Delta_1 v(t = 5 \text{ ms})} \right| \quad (36)$$

in which $\Delta_1 v$ is the velocity dispersion extracted by the full numerical approach, and $\Delta_2 v$ is the one resulting from the rescaled initial Wigner function. Results are summarized in the following table:

N	176	880	1760	8800	17600	35200
ε	0.278	0.0741	0.0397	0.0132	0.0104	0.009

Table 1: Numerical values of $\varepsilon(N)$.

The larger the interaction parameter χ ($\propto N$), the better the agreement. We conclude that our scaling approach gives a good account of the evolution of the wave function of a Bose-Einstein condensate in the Thomas-Fermi regime. As the validity of our approach is now well demonstrated, we can extend its use for much larger time. We introduce in the following the quantity \mathcal{M}_x which characterized the improvement of the local monochromaticity as a function of the distance from the original trap:

$$\mathcal{M}_x = \frac{\Delta v_x}{\langle v \rangle_x} = \frac{\sqrt{\int_0^\infty V_3(x, t) dt - \left(\int_0^\infty V_2(x, t) dt \right)^2}}{\int_0^\infty V_2(x, t) dt} \quad (37)$$

To calculate the moments involved in the local monochromaticity quantity, \mathcal{M}_x , we use the scaling laws provided

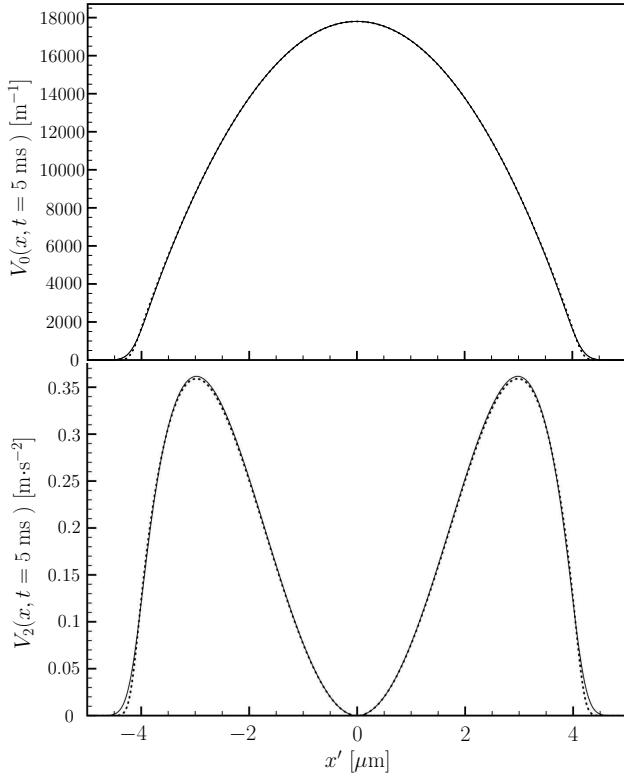


FIG. 5: Velocity moments $V_0(x, t = 5 \text{ ms})$ (density profile) and $V_2(x, t = 5 \text{ ms})$ for $N = 35200$ atoms, where the positions are expressed in the center-of-mass coordinates $x' = x - \eta$; full numerical procedure including the time evolution (solid line), result of the scaling approach based only on the initial Wigner function (dotted line).

by Eqs. (21) and (22). The evolution of this quantity as a function of the distance is plotted in Fig. 6. We recover here quantitatively the main result of this section, the local monochromaticity increases with the distance even for a many-body wave function with large repulsive interaction. This is to be contrasted with the global velocity dispersion which increases as shown in Fig. 4. An interacting BEC evolving in the presence of a constant force can therefore be used as a local probe in a scattering experiment. In the following section, we give a concrete example in which the outcoupling of the wave function from its original trap is performed progressively to amplify the local monochromaticity including in the case of a Bose-Einstein condensate in the Thomas-Fermi regime.

VI. GAIN OBTAINED FROM A PROGRESSIVE OUTCOUPLING

In this section, we compare the matter wave probing of a very thin Fabry-Perot resonance associated with a repulsive double barrier [21, 22] using two different protocols: first we consider a wave packet at a constant ve-

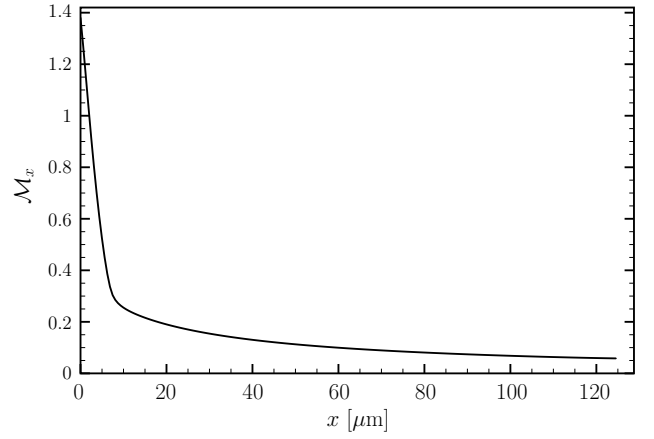


FIG. 6: Local monochromaticity as a function of the distance x at which it is measured. This data has been obtained for a Bose-Einstein condensate in the Thomas-Fermi regime with $N = 35200$ atoms, other parameters are identical to those of Fig. 3.

locity with a well-defined velocity dispersion and in the absence of other extra external potential (see Fig. 7.a); second we consider a situation in which the matter wave is accelerated by a linear potential towards the double barrier (see Fig. 8.b) and for which atoms can also be progressively outcoupled from their original trap [23–27].

The double barrier potential is modeled by a sinusoidal potential located at $x = d$:

$$U_{\text{db}}(x) = 2sE_R(1 - \cos(2\pi(x - d)/d_R)) \times \Theta(x - d + d_R)\Theta(d + d_R - x), \quad (38)$$

where $E_R = \hbar^2/(2md_R^2) = mv_R^2/2$, Θ is the Heaviside step function ($\Theta(x < 0) = 0$ and $\Theta(x \geq 0) = 1$). The numerical simulations presented hereafter are performed with the realistic values $d = 201 \text{ } \mu\text{m}$, $d_R = 0.48 \text{ } \mu\text{m}$, $v_R = \hbar/(md_R) = 9.565 \text{ mm/s}$ [7, 28]. The Fabry-Perot resonance that is probed has an energy $E_0 = mv_0^2/2$ with $v_0 = 1.62v_R = 15.53 \text{ mm/s}$ for a barrier height of $16.08E_R = 6.13E_0$. The width of this resonance is $E_0/1000$ (i.e. 0.36 pK in temperature units!) that corresponds to a velocity dispersion $8.3 \text{ } \mu\text{m/s}$. For comparison, let us recall that the typical order of magnitude of the velocity dispersion achievable with a rubidium BEC is 2 mm/s [7].

To scan the resonance we vary the height of the double barrier by adjusting the parameter s for a fixed incoming velocity. The result obtained by considering an incident plane wave of wave vector $k_0 = mv_0/\hbar$ is plotted as a reference by a solid line in Figs. 7(b) and 8(b). The resonance appears at $s = 4.02$. Interestingly, the width in energy of the reflection probability as a function of s is significantly larger than that of the resonance. Indeed, from a Gaussian fit of the reflection probability, we find a variance in s equal to $\Delta s = 0.008$. This means that experimentally if the double barrier is realized by optical

means, the intensity should be stabilized at a level better than one per thousand, and not much better than the resonance width $E_0/1000$. This feature facilitates the detection of the resonance.

Figure 7(b) contains also the reflection probability about the resonance for an incident wave packet of mean velocity v_0 and with a velocity dispersion $\Delta v = 600 \mu\text{m/s}$ (dashed line) and $20 \mu\text{m/s}$ (dotted line). The probability reflection of an incident packet with a velocity dispersion already more than three times smaller than that of a BEC remains very close to one. This is to be contrasted with the case of a packet that have a velocity dispersion 100 times smaller than that of a BEC (i.e. on the order of the width of the resonance). In this latter case, the presence of the resonance can be clearly seen as expected since about 38 % of the atoms are reflected.

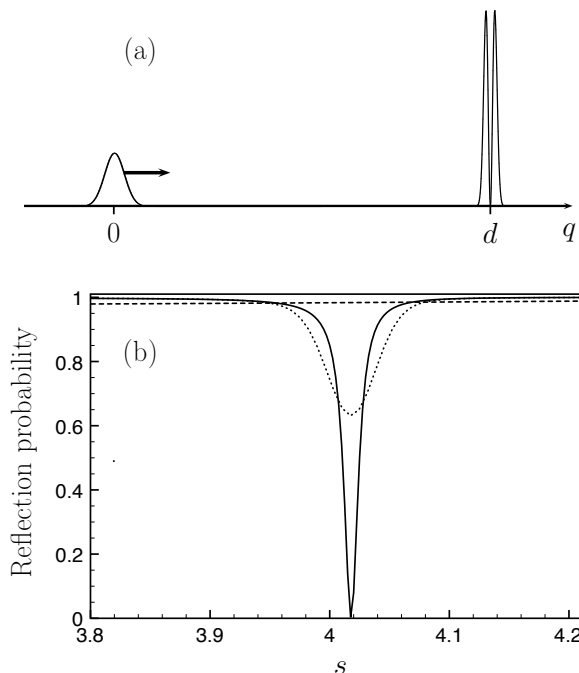


FIG. 7: Matter wave probing of a Fabry-Perot resonance. The reflection probability is plotted as a function of the double barrier height: for an incoming plane wave (solid line) with the velocity v_0 that coincides with the resonance ($s = 4.02$), for an incident wave packet with a velocity dispersion of $600 \mu\text{m/s}$ (dashed line) and $20 \mu\text{m/s}$ (dotted line).

We now consider the probing of the same resonance by a wave packet that is accelerated by a constant force from its original trap located at $x = 0$. The potential energy experienced by the atoms now reads (see Fig. 8(a))

$$U(x, t) = m\gamma(x - d) - U_0 e^{-t/\tau} e^{-2x^2/w_0^2} + U_{\text{db}}(x). \quad (39)$$

The Gaussian term accounts for the trap potential which accommodates the initial wave function. The exponential form for the time dependence of the trap depth was used in Ref. [29] to outcouple progressively atoms from a Bose-Einstein condensate into a guide. The slope has been

chosen so that a particle launched with a zero velocity at $x = 0$ arrives at $x = d$ with a velocity v_0 i.e. $\gamma = v_0^2/(2d) = 0.6 \text{ m/s}^2$ i.e. an energy that coincides with that of the Fabry-Perot resonance.

The analytical formalism developed in the previous sections is the appropriate one for $\tau \rightarrow 0$, i.e. for an instantaneous release of the wave function from its originating trap and a subsequent propagation on the slope. We have already shown that the position-velocity correlation that builds up during such an evolution generates a small local velocity dispersion downward. As demonstrated below on a specific example, an even better strategy consists in releasing the atoms from the trap progressively as done for the generation of guided atom lasers [28–32]. In such a protocol, the initial conditions of each atom is nearly the same i.e. atoms are released nearly at the same position with a velocity very close to zero. This favors a very thin local velocity dispersion after propagation.

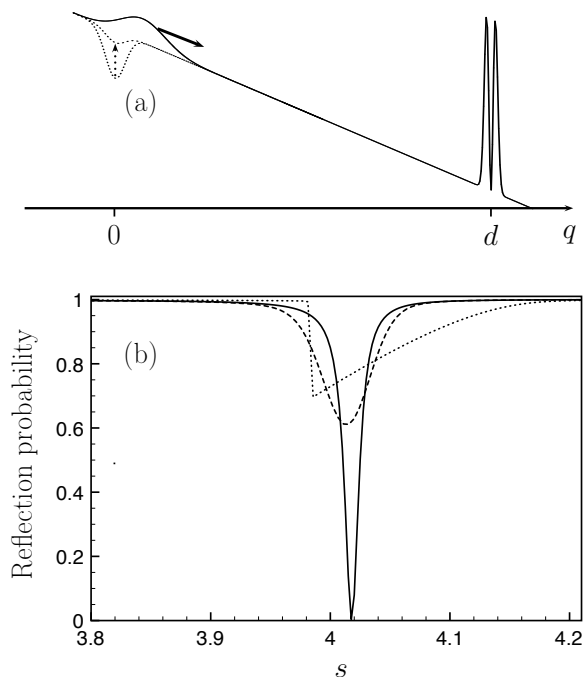


FIG. 8: Matter wave probing of a Fabry-Perot resonance placed on a slope. The reflection probability is plotted as a function of the double barrier height: for an incoming plane wave (solid line) with the velocity v_0 that coincides with the resonance ($s = 4.02$), for an incident wave packet with a velocity dispersion of $600 \mu\text{m/s}$ outcoupled in 80 ms (dashed line) and for a Bose-Einstein condensate in the Thomas-Fermi of velocity dispersion 2 mm/s outcoupled in 100 ms (dotted line).

We illustrate this idea with the dashed curve of Fig. 8.b that corresponds to the reflection probability of a wave packet (without atom-atom interactions) with an initial velocity dispersion equal to $600 \mu\text{m/s}$ and that has been outcoupled in $\tau = 80 \text{ ms}$. About 40 % of the atoms are

reflected when the height of the double barrier is varied. In contrast with the scattering on the double barrier, in the absence of slope, of a packet having the same velocity dispersion, the thin Fabry resonance can here be probed with accuracy. This illustrates clearly the importance of both a progressive outcoupling and the building up of position-velocity correlation in the course of the propagation on the slope.

We have also performed the progressive outcoupling (over 100 ms) of an interacting BEC with a velocity dispersion even larger on the order of 2 mm/s [18]. Note that the long outcoupling times are also optimal to reduce the contamination of the transverse excited states in real systems [18, 28, 29, 31]. The result is plotted as a dotted line in Fig. 8(b). The reflection probability is about 30 %. Once again the initial source is very far from an ideal one and has a “very” large instantaneous velocity width. The situation may appear as even worse since the interaction energy is converted into kinetic energy in the course of the propagation as already discussed in Sec. V. However, the benefit of the building up of the position-velocity correlation is dominant. An instantaneous outcoupling of the BEC yields a 10 % reflection probability, a value that would correspond in the absence of a slope to that of a wave packet having a dispersion velocity on the order of 50 $\mu\text{m/s}$, that is 40 times smaller than that of the BEC. An extra gain of a factor three on the reflection probability is here obtained from the progressive outcoupling. Interestingly, this technique avoids completely any limitation in energy dictated by the chemical potential. The shape of the reflection probability as a function of the height of the double barrier is radically different in the case of an interacting BEC compared to that obtained in the absence of atom-atom interactions (compare dotted and dashed lines in Fig. 8.b). This results from the sharp border of the inverse parabolic shape of the BEC assumed in the Thomas-Fermi regime and its conversion in position-velocity correlation in the course of the propagation.

In conclusion, we have shown how it is possible to decrease the velocity dispersion at a given position using a linear potential. This method works efficiently also in the presence of repulsive interaction. It provides a method to circumvent the limitations that may appear in outcoupling processes [22, 34] and to improve the local monochromaticity of a matter wave source.

In Ref. [33], the author proposed to investigate an atom-blockade effect with a Fabry-Perot cavity in close analogy with the Coulomb blockade effect of electronic transport. An incident atom laser with a very thin velocity dispersion would be reflected or not from a Fabry-Perot cavity depending on the presence or not of an atom in the cavity. Such an experiment appeared as out of reach for a long time because of the very small velocity dispersion ($< 50 \mu\text{m/s}$) required for its demonstration. Exploiting the position-velocity correlation that builds up in the propagation over a slope we have shown that this level of accuracy can now be achieved with the cur-

rent state of the art. This work therefore paves the way for ultrahigh resolution in matter wave probing, the observation of nonlinear atom optical effects and of the interplay between atom tunneling and atom-atom interactions [35, 36].

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Appendix A: Scaling Wigner solution

This appendix justifies the scaling form of the Wigner function of Eq. (15) for both a one-body wave function and a Bose-Einstein condensate in the Thomas-Fermi regime.

We consider a cloud of atoms in a harmonic trap. At time $t = 0$, the confinement is suddenly removed and the packet experiences a uniform accelerating force. For a non-interacting gas, the initial Wigner function, $W_0(x, p)$, obeys the stationary Liouville equation

$$p\partial_x W_0 - m^2\omega_0^2 x\partial_p W_0 = 0, \quad (\text{A1})$$

Multiplying this by p/N , where N is the total number of particles, and integrating over p yields

$$\partial_x \left(\int p^2 W_0 dp \right) + m^2\omega_0^2 x n_0(x) = 0, \quad (\text{A2})$$

where $n_0(x)$ is the linear density defined by

$$n_0(x) = N \int W_0(x, p) dp \quad \text{with} \quad \int n_0(x) dx = N. \quad (\text{A3})$$

Multiplying Eq. (A2) by x/N and integrating over x provides the usual momentum relation associated to the harmonic oscillator $\langle p^2 \rangle_0 = m^2\omega_0^2 \langle x^2 \rangle_0$.

For $t > 0$, the packet is released on a slope i.e. on a linear potential $U(x) = -m\gamma x$. The equation of evolution of the time-dependent Wigner function now reads

$$\partial_t W + (p/m)\partial_x W + m\gamma\partial_p W = 0. \quad (\text{A4})$$

We look for the solution of this time-dependent equation using a scaling ansatz : $W(x, p, t) = W_0(X, P)$ with $X = [x - \eta]/\alpha$ and $P = \alpha(p - m\dot{\eta}) - m\dot{\alpha}(x - \eta)$ [10], where α and η are time-dependent functions to be determined with boundary conditions

$$\eta(0) = 0, \quad \dot{\eta}(0) = 0, \quad \alpha(0) = 1, \quad \dot{\alpha}(0) = 0, \quad (\text{A5})$$

so that $P(t = 0) = p$ and $X(t = 0) = x$. Let’s calculate the derivative of X and P variables with respect to x, p

and t :

$$\begin{aligned}
\partial_t X &= -(\dot{\alpha}X + \dot{\eta})/\alpha \\
\partial_x X &= 1/\alpha \\
\partial_t P &= m\dot{\alpha}\dot{\eta} - m\alpha\ddot{\eta} + mX(\dot{\alpha}^2 - \alpha\ddot{\alpha}) + \dot{\alpha}P/\alpha \\
\partial_x P &= -m\dot{\alpha} \\
\partial_P P &= \alpha.
\end{aligned} \tag{A6}$$

We deduce the following derivative of the phase-space distribution :

$$\begin{aligned}
\partial_t W &= \partial_t X \partial_X W_0 + \partial_t P \partial_P W_0, \\
\partial_x W &= \partial_x X \partial_X W_0 + \partial_x P \partial_P W_0, \\
\partial_P W &= \partial_P P \partial_P W_0.
\end{aligned} \tag{A7}$$

Combining Eqs. (A7) and (A4), we get

$$\Lambda_X \partial_X W_0 + \Lambda_P \partial_P W_0 = 0 \tag{A8}$$

with

$$\begin{aligned}
\Lambda_X &= (\partial_t X + (p/m)\partial_x X) = P/(m\alpha^2), \\
\Lambda_P &= (\partial_t P + (p/m)\partial_x P + m\gamma\partial_P P) \\
&= -m\alpha(\ddot{\eta} - \gamma) - mX\alpha\ddot{\alpha}.
\end{aligned} \tag{A9}$$

The integration of Eq. (A8) over X and P yields the equation for the packet center $\ddot{\eta} = \gamma$. Multiplying Eq. (A8) by XP and integrating over X and P provides the equation fulfilled by the α scaling parameter:

$$\ddot{\alpha} = \frac{\langle P^2 \rangle_0}{m^2 \langle X^2 \rangle_0} \frac{1}{\alpha^3} \equiv \frac{\omega_0^2}{\alpha^3}, \tag{A10}$$

whose solution is $\alpha(t) = (1 + \omega_0^2 t^2)^{1/2}$. For the ground state we find $\omega_0 = \hbar/(2m\sigma^2)$ where $\sigma = \Delta x_0$.

Let us now consider a Bose-Einstein condensate in the Thomas-Fermi regime. Initially the BEC is trapped in an harmonic trap of angular frequency ω_0 and its wave function normalized to the number of atoms reads $\Psi(x, 0) = n_0^{1/2}(x)$, where n_0 is the atomic density:

$$n_0(x) = (\mu - m\omega_0^2 x^2/2)/g, \tag{A11}$$

where g accounts for the strength of the interactions. At time $t > 0$, we propose to describe the dynamics using the ansatz

$$\Psi(x, t) = \frac{1}{\alpha^{1/2}} n_0^{1/2} \left(\frac{x - \eta}{\alpha} \right) e^{iS(x, t)}. \tag{A12}$$

To obtain the expression of the scaling parameter α and of the phase $S(x, t)$ we shall use the hydrodynamic equations. The continuity equation for the density $n(x, t) = |\Psi(x, t)|^2$ yields the expression for the velocity field [11],

$$\partial_t n + \partial_x (nv_x) = 0 \implies v_x = \dot{\eta} + \dot{\alpha} \frac{x - \eta}{\alpha}. \tag{A13}$$

The phase S can be directly inferred from the velocity field through the relation $v_x = \hbar \partial_x S / m$:

$$S(x, t) = \frac{m}{\hbar} \left(\dot{\eta}(x - \eta) + \dot{\alpha} \frac{(x - \eta)^2}{2\alpha} \right). \tag{A14}$$

The Euler equation for the velocity field in the Thomas Fermi regime reads

$$m\partial_t v_x = \partial_x \left(-\frac{1}{2}mv^2 - V_{\text{ext}}(x) - gn(x, t) \right). \tag{A15}$$

Combining Eqs. (A11), (A12) and (A15), we find the equations fulfilled by the center position of the packet η and the one for scaling parameter α ,

$$\ddot{\eta} = \gamma \quad \text{and} \quad \ddot{\alpha} = \frac{\omega_0^2}{\alpha^2}. \tag{A16}$$

Instead of the α^3 denominator that we found for the noninteracting gas in Eq. (A10), we have now α^2 , so the solution cannot be written explicitly. However, as $\dot{\alpha} = \omega_0[2(1 - \alpha^{-n+1})/(n - 1)]^{1/2}$, the long time asymptotics, as $t \gg 1/\omega_0$, is quite simple: $\dot{\alpha} \sim \omega_0$ for the non-interacting gas ($n = 3$) and $\dot{\alpha} \sim \sqrt{2}\omega_0$ for the TF regime ($n = 2$). Once the mean field energy is released, the cloud expansion proceeds as a free expansion with a rate of change corresponding to an effective initial frequency $\sqrt{2}\omega_0$, a fingerprint of the interactions that, remarkably, does not depend on g .

The time-dependent Wigner function associated with $\Psi(x, t)$ can be recast in terms of the initial Wigner function:

$$\begin{aligned}
W(x, p, t) &= \frac{1}{N\pi\hbar} \int \Psi^*(x + y, t) \Psi(x - y, t) e^{2ipy/\hbar} dy \\
&= \frac{1}{N\pi\hbar\alpha} \int \left[n_0 \left(\frac{x + y - \eta}{\alpha} \right) n_0 \left(\frac{x - y - \eta}{\alpha} \right) \right]^{1/2} \\
&\quad \times e^{2ipy/\hbar + iS(x - y, t) - iS(x + y, t)} dy \\
&= W_0(X, P).
\end{aligned} \tag{A17}$$

Interestingly enough, this result coincides with the one of the collisionless Vlasov equation [37]. This may appear surprising at first sight since Vlasov equation is a semi-classical equation in the presence of a mean-field term. However, the quadratic form of the density profile that enters the mean field potential term enables to conserve the equivalence between the equation of evolution for the Wigner function and the Vlasov equation. In this sense, this is a direct consequence of the initial harmonic confinement assumption.

The expression used above for the Wigner is the one associated with a pure state. This is valid in the absence of interactions and also for the mean field description of a Bose-Einstein condensate but it cannot be applied directly to a many-body wave function. In this latter case, the Wigner function is defined through the one-body reduced density matrix, $g_1(x, y; t)$ [9]:

$$W(x, p, t) = \frac{1}{\pi\hbar} \int g_1(x + y, x - y, t) e^{2ipy/\hbar} dy. \tag{A18}$$

A large class of many-body quantum systems exhibit self-similar dynamics [13, 38]. This includes the Calogero-Sutherland model [39], the Tonks Girardeau gas [40, 41],

certain Lieb-Liniger states [42], Bose-Einstein condensate [43, 44] even in the presence of dipolar interactions [45], strongly interacting gas mixtures [46], ... The self-similar dynamics can be written in our context as

$$g_1(x, y; t) = \frac{1}{\alpha} g_1\left(\frac{x - \eta}{\alpha}, \frac{y - \eta}{\alpha}; 0\right) e^{i(S(x, t) - S(y, t))} \quad (\text{A19})$$

yielding once again $W(x, p, t) = W_0(X, P)$.

Appendix B: f functions

We may calculate the f_j functions in Eq. (30) making use of Eqs. (20-25). An alternative route is to write down the Wigner function explicitly making use of the known phase-space propagator,

$$\begin{aligned} W(x, p; t) &= \iint dx_0 dp_0 W(x_0, p_0; 0) \\ &\times \delta\left(x - x_0 - v_0 t - \frac{\gamma t^2}{2}\right) \delta(v - v_0 - \gamma t) \\ &= \frac{1}{\pi \hbar} e^{-\frac{(p - m\gamma t)^2 2\sigma^2}{\hbar^2}} e^{-\frac{(x - pt/m + \gamma t^2/2)^2}{2\sigma^2}}, \quad (\text{B1}) \end{aligned}$$

$$f_1 = \frac{m\sigma t(8\gamma m^2 \sigma^4 + \gamma \hbar^2 t^2 + 2\hbar^2 x)}{\sqrt{2\pi}(4m^2 \sigma^4 + \hbar^2 t^2)^{3/2}}, \quad (\text{B2})$$

$$f_2 = \frac{m\sigma \{64\gamma^2 m^4 \sigma^8 t^2 + 16\hbar^2 m^2 \sigma^4 [\sigma^2 + \gamma t^2(\gamma t^2 + 2x)] + \hbar^4 t^2 [4\sigma^2 + (\gamma t^2 + 2x)^2]\}}{2\sqrt{2\pi}(4m^2 \sigma^4 + \hbar^2 t^2)^{5/2}}, \quad (\text{B3})$$

$$\begin{aligned} f_3 &= \frac{m\sigma t(8\gamma m^2 \sigma^4 + \gamma \hbar^2 t^2 + 2\hbar^2 x)}{4\sqrt{2\pi}(4m^2 \sigma^4 + \hbar^2 t^2)^{7/2}} \\ &\times \left\{ 64\gamma^2 m^4 \sigma^8 t^2 + 16\hbar^2 m^2 \sigma^4 [3\sigma^2 + \gamma t^2(\gamma t^2 + 2x)] + \hbar^4 t^2 [12\sigma^2 + (\gamma t^2 + 2x)^2] \right\}. \quad (\text{B4}) \end{aligned}$$

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